LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
Data File : VX030276.D
Acq On : 28 Jul 2022 17:45
Operator : JC/MD
Sample : N3953-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Title : SW846 8260

Signal : TIC: VX030276.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.142	6	9	16	rVB3	9775	11383	1.03%	0.098%
2	1.239	22	25	27	rBV	19039	22788	2.06%	0.196%
3	1.355	40	44	52	rVB2	12394	16434	1.49%	0.141%
4	1.453	56	60	69	rVB	246427	280493	25.37%	2.414%
5	2.063	154	160	170	rBV	206351	313511	28.36%	2.698%
6	2.294	193	198	207	rVB	47747	75000	6.78%	0.645%
7	2.386	207	213	225	rBV	191911	318110	28.78%	2.738%
8	2.544	228	239	245	rBV	27978	54170	4.90%	0.466%
9	2.709	259	266	275	rVB2	8626	16222	1.47%	0.140%
10	2.971	301	309	318	rVB2	4811	12522	1.13%	0.108%
11	3.117	320	333	343	rBV	141194	323530	29.27%	2.784%
12	3.331	362	368	379	rVB4	5714	13198	1.19%	0.114%
13	3.581	402	409	419	rBV4	9434	23394	2.12%	0.201%
14	3.867	450	456	472	rBV4	9074	26876	2.43%	0.231%
15	4.245	507	518	522	rBV3	13330	35843	3.24%	0.308%
16	4.294	522	526	532	rVV	13364	33177	3.00%	0.286%
17	4.367	532	538	549	rVB4	19447	57346	5.19%	0.493%
18	4.568	562	571	587	rBV2	47275	134558	12.17%	1.158%
19	5.391	695	706	720	rBV	136498	402572	36.42%	3.464%
20	5.562	724	734	746	rVV2	143773	415791	37.61%	3.578%
21	5.964	790	800	811	rBV	165558	445792	40.33%	3.836%
22	6.647	904	912	920	rBV	9401	23442	2.12%	0.202%
23	6.769	921	932	943	rVV	240826	582382	52.68%	5.012%
24	7.159	989	996	1003	rVB2	7424	16491	1.49%	0.142%
25	7.470	1041	1047	1054	rVV	8574	19032	1.72%	0.164%
26	7.586	1059	1066	1073	rBV2	69742	140126	12.68%	1.206%
27	8.580	1222	1229	1234	rBV2	14759	23801	2.15%	0.205%
28	8.653	1234	1241	1248	rVV	732719	1105461	100.00%	9.513%
29	8.720	1248	1252	1260	rVB	41137	70877	6.41%	0.610%
30	9.037	1301	1304	1311	rVB2	8496	15364	1.39%	0.132%
31	9.244	1333	1338	1343	rBV4	9178	17136	1.55%	0.147%
32	9.525	1377	1384	1398	rBV	120188	177736	16.08%	1.530%
33	10.055	1465	1471	1482	rVB	514171	696853	63.04%	5.997%
34	10.201	1490	1495	1504	rBV2	13281	24883	2.25%	0.214%
35	10.305	1504	1512	1519	rVB	56316	82434	7.46%	0.709%
36	10.591	1554	1559	1563	rBV3	12429	20881	1.89%	0.180%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Title : SW846 8260

37	10.646	1563	1568	1572	rVV	40026	51229	4.63%	0.441%
38	10.774	1584	1589	1592	rBV	10660	15381	1.39%	0.132%
39	10.860	1598	1603	1608	rBV4	12459	19938	1.80%	0.172%
40	11.085	1634	1640	1647	rBV	534597	658010	59.52%	5.663%
41	11.384	1684	1689	1691	rBV	29140	42150	3.81%	0.363%
42	11.457	1697	1701	1710	rVB	25969	38538	3.49%	0.332%
43	11.616	1721	1727	1732	rBV2	40850	59293	5.36%	0.510%
44	11.756	1743	1750	1756	rVB	108551	146982	13.30%	1.265%
45	12.024	1789	1794	1800	rBV	548984	654250	59.18%	5.630%
46	12.091	1800	1805	1811	rVB	51881	67331	6.09%	0.579%
47	12.244	1823	1830	1833	rBV4	50163	90735	8.21%	0.781%
48	12.317	1838	1842	1850	rVB3	36677	55958	5.06%	0.482%
49	12.427	1854	1860	1863	rBV5	13802	19410	1.76%	0.167%
50	12.463	1863	1866	1872	rVB	11694	15042	1.36%	0.129%
51	12.536	1873	1878	1879	rBV	29108	34790	3.15%	0.299%
52	12.555	1879	1881	1885	rVB	20464	23992	2.17%	0.206%
53	12.616	1885	1891	1895	rBV	57393	74201	6.71%	0.639%
54	12.701	1901	1905	1913	rVB2	23654	44220	4.00%	0.381%
55	12.811	1913	1923	1926	rBV8	17719	44122	3.99%	0.380%
56	12.853	1926	1930	1934	rVB2	26604	37644	3.41%	0.324%
57	12.926	1936	1942	1945	rVV	54949	72800	6.59%	0.626%
58	12.963	1945	1948	1953	rVB	96971	123478	11.17%	1.063%
59	13.048	1955	1962	1964	rBV5	13760	25201	2.28%	0.217%
60	13.170	1978	1982	1987	rVB	59029	76107	6.88%	0.655%
61	13.292	1992	2002	2007	rVV3	140757	262807	23.77%	2.262%
62	13.347	2007	2011	2020	rVB4	66922	106712	9.65%	0.918%
63	13.426	2020	2024	2028	rBV3	17489	24314	2.20%	0.209%
64	13.506	2033	2037	2040	rBV4	16297	26302	2.38%	0.226%
65	13.548	2040	2044	2046	rVV	35388	43612	3.95%	0.375%
66	13.585	2046	2050	2057	rVV5	40270	87846	7.95%	0.756%
67	13.670	2057	2064	2068	rVV3	27588	61433	5.56%	0.529%
68	13.713	2068	2071	2075	rVV2	25285	32859	2.97%	0.283%
69	13.780	2077	2082	2088	rVV	355201	463875	41.96%	3.992%
70	13.835	2088	2091	2100	rVB5	19367	42074	3.81%	0.362%
71	13.926	2102	2106	2110	rBV4	24682	36831	3.33%	0.317%
72	13.975	2110	2114	2117	rVV2	32255	39916	3.61%	0.344%
73	14.042	2121	2125	2127	rVV	31479	30364	2.75%	0.261%
74	14.079	2127	2131	2134	rVB	42654	46060	4.17%	0.396%
75	14.182	2143	2148	2150	rBV2	23806	31079	2.81%	0.267%
76	14.213	2150	2153	2158	rVV	43185	68015	6.15%	0.585%

LSC Area Percent Report

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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Title : SW846 8260

77	14.262	2158	2161	2167	rVB2	34507	48513	4.39%	0.417%
78	14.323	2167	2171	2175	rBV	37132	44120	3.99%	0.380%
79	14.371	2176	2179	2183	rBV	33885	43287	3.92%	0.373%
80	14.414	2183	2186	2192	rVB6	13264	25366	2.29%	0.218%
81	14.481	2193	2197	2200	rBV6	15192	24298	2.20%	0.209%
82	14.640	2216	2223	2228	rBV	524016	640752	57.96%	5.514%
83	14.780	2234	2246	2254	rBV	327146	462055	41.80%	3.976%
84	14.883	2260	2263	2269	rVB5	9253	17243	1.56%	0.148%
85	14.993	2277	2281	2286	rVB8	8724	13761	1.24%	0.118%
86	15.213	2314	2317	2323	rVB5	14740	18917	1.71%	0.163%
87	15.377	2339	2344	2347	rBV3	21431	34996	3.17%	0.301%
88	15.481	2355	2361	2370	rBV4	47819	94057	8.51%	0.809%
89	15.615	2378	2383	2387	rBV2	52342	87806	7.94%	0.756%
90	15.652	2387	2389	2396	rVB2	28861	39485	3.57%	0.340%
91	15.841	2412	2420	2425	rBV3	15679	37106	3.36%	0.319%
92	15.987	2439	2444	2450	rBV3	15124	28838	2.61%	0.248%
93	16.090	2457	2461	2470	rBV10	7094	15252	1.38%	0.131%

Sum of corrected areas: 11620362

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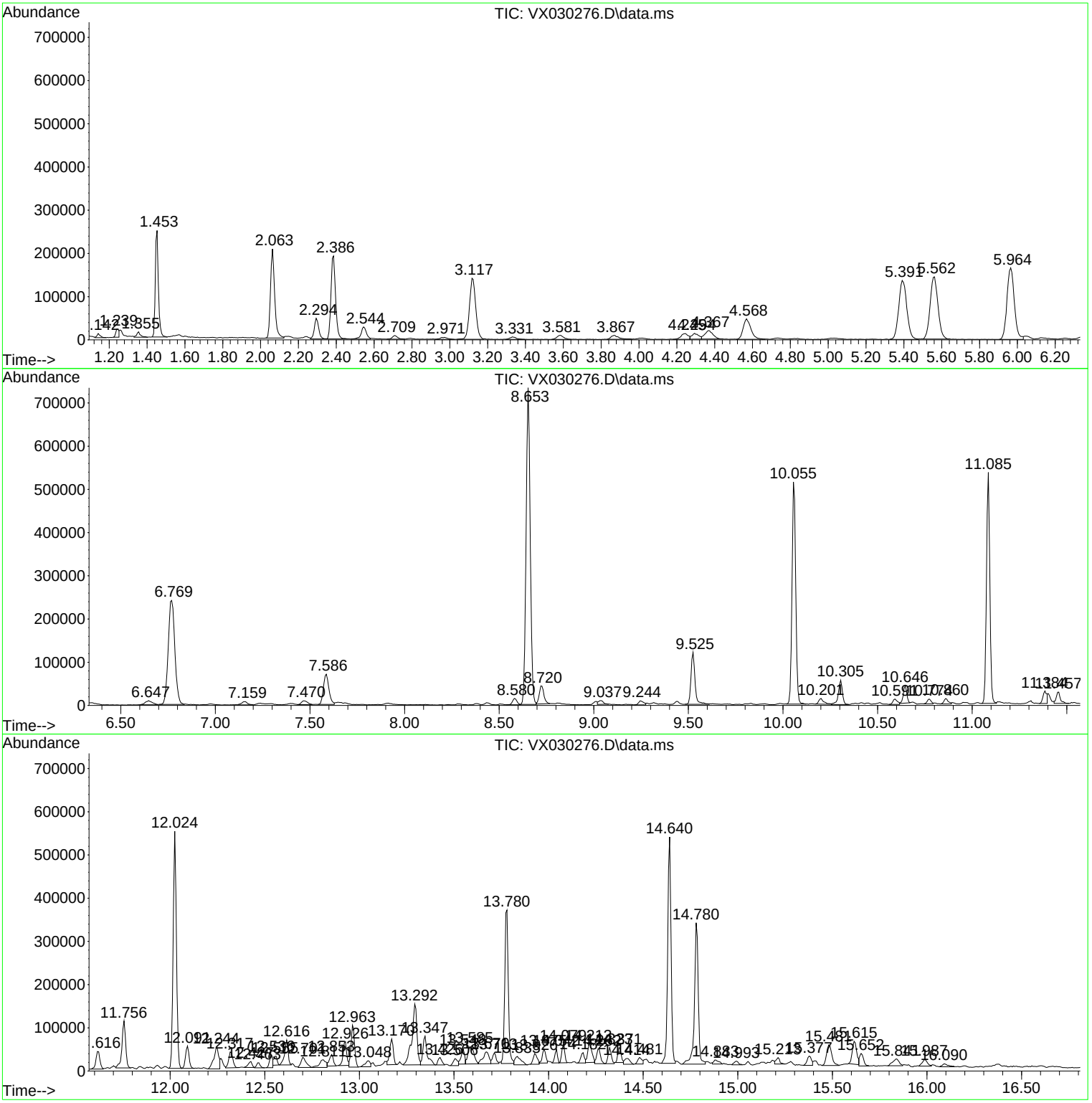
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
Data File : VX030276.D
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Operator : JC/MD
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ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

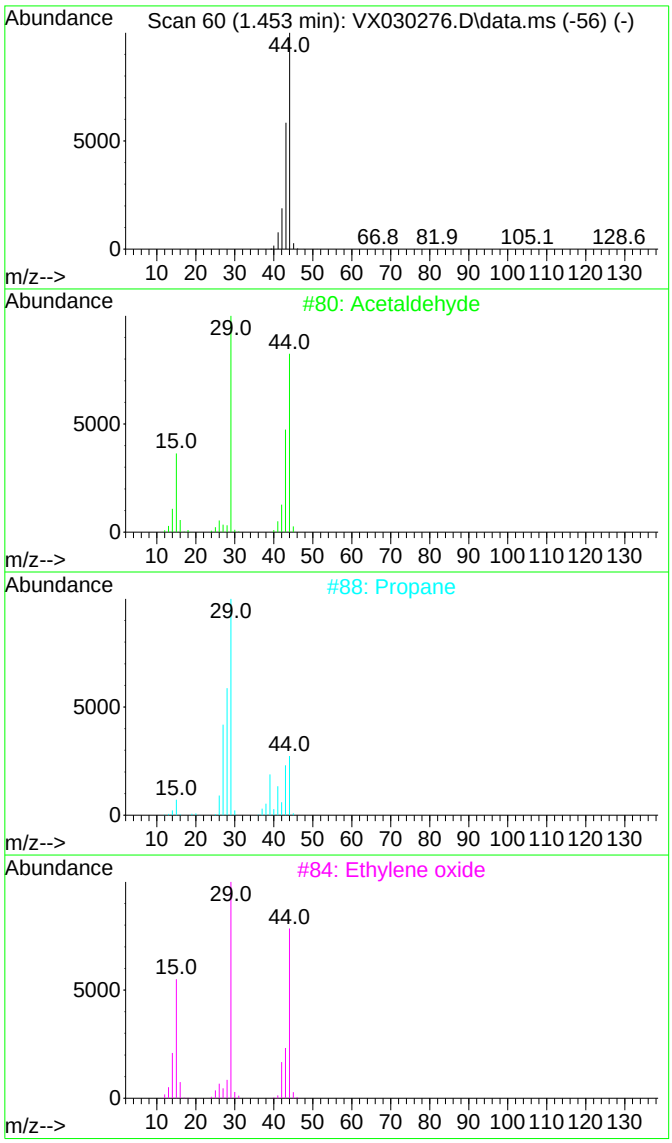
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

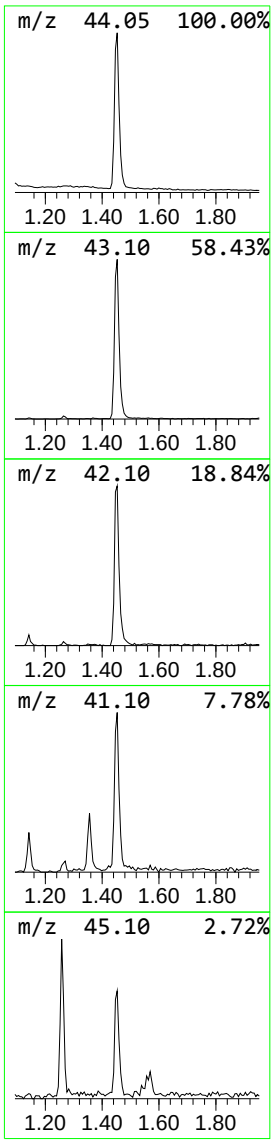
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Acetaldehyde Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.453	33.73 ug/l	280493	Pentafluorobenzene	5.562

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetaldehyde	44	C2H4O	000075-07-0	64
2	Propane	44	C3H8	000074-98-6	5
3	Ethylene oxide	44	C2H4O	000075-21-8	5
4	Alanine	89	C3H7NO2	000056-41-7	4
5	Hydrogen isocyanate	43	CHNO	000075-13-8	3





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
Data File : VX030276.D
Acq On : 28 Jul 2022 17:45
Operator : JC/MD
Sample : N3953-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

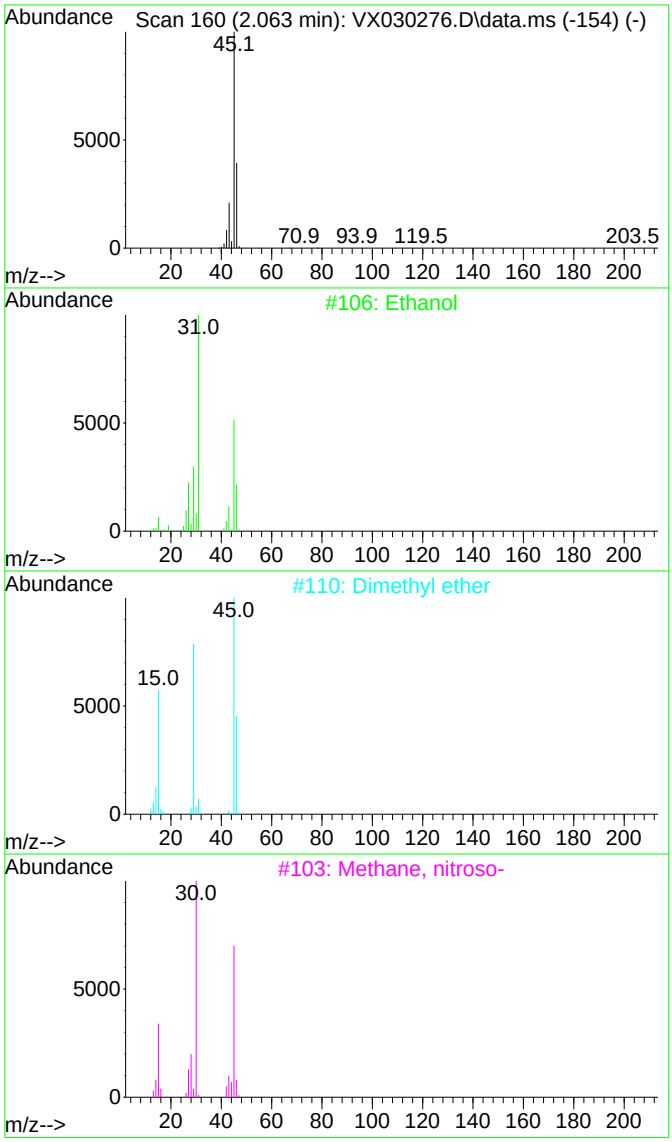
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

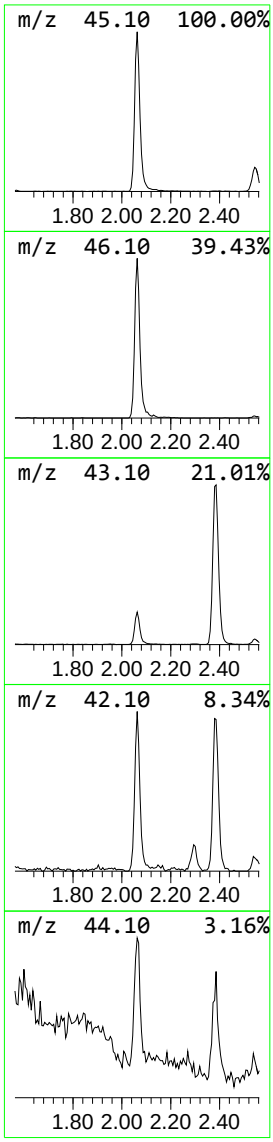
****

Peak Number 2 Ethanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.063	37.70 ug/l	313511	Pentafluorobenzene	5.562

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol	46	C2H6O	000064-17-5	64
2	Dimethyl ether	46	C2H6O	000115-10-6	9
3	Methane, nitroso-	45	CH3NO	000865-40-7	4
4	Formic acid	46	CH2O2	000064-18-6	4
5	Oxalic acid	90	C2H2O4	000144-62-7	4





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
Data File : VX030276.D
Acq On : 28 Jul 2022 17:45
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

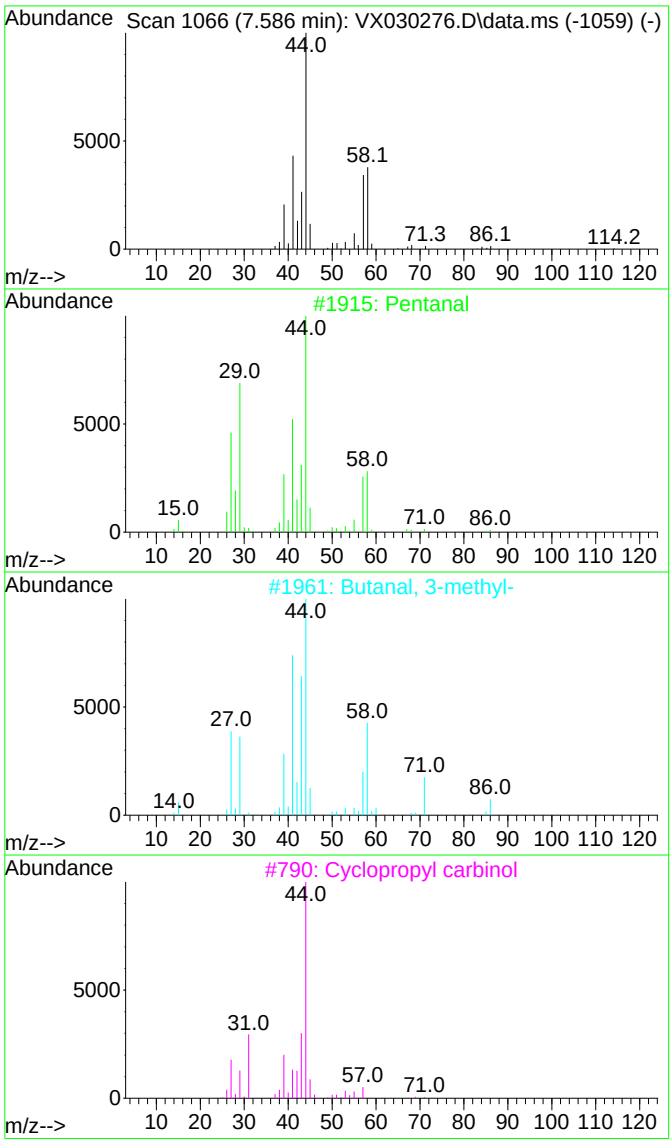
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

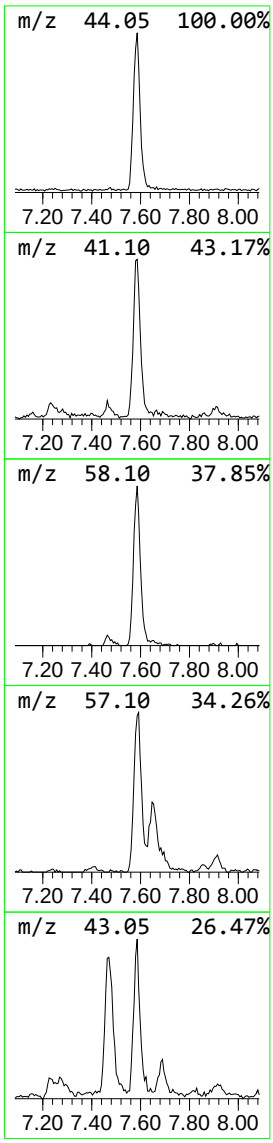
****

Peak Number 3 Pentanal Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.586	12.03 ug/l	140126	1,4-Difluorobenzene	6.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanal	86	C5H10O	000110-62-3	90
2	Butanal, 3-methyl-	86	C5H10O	000590-86-3	74
3	Cyclopropyl carbinol	72	C4H8O	002516-33-8	59
4	Piperazine	86	C4H10N2	000110-85-0	38
5	2-Formylhistamine	139	C6H9N3O	1000132-95-8	36





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
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Acq On : 28 Jul 2022 17:45
Operator : JC/MD
Sample : N3953-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

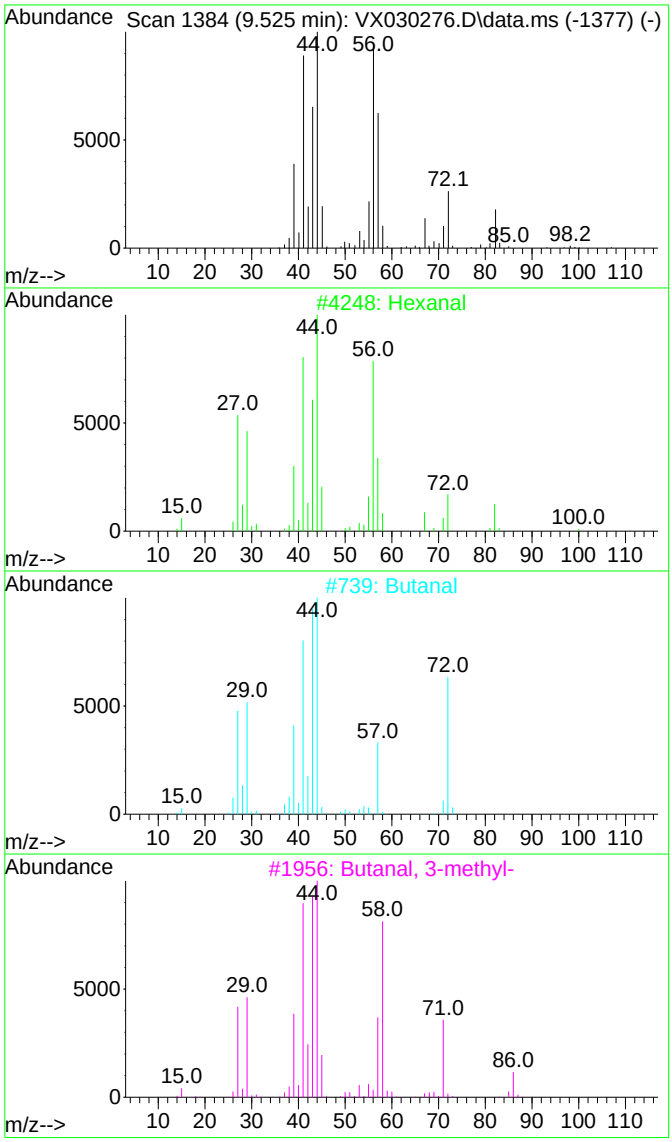
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

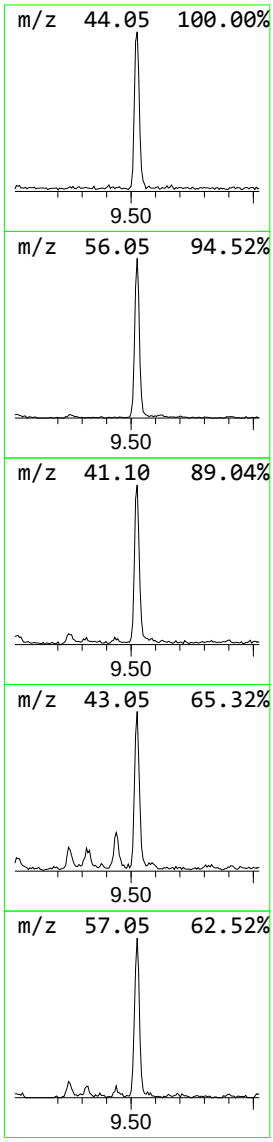
****

Peak Number 4 Hexanal Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.525	12.75 ug/l	177736	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal	100	C6H12O	000066-25-1	72
2	Butanal	72	C4H8O	000123-72-8	43
3	Butanal, 3-methyl-	86	C5H10O	000590-86-3	37
4	Cyclopentanol, 2-methyl-, cis-	100	C6H12O	025144-05-2	35
5	Glutaraldehyde	100	C5H8O2	000111-30-8	27





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Data File : VX030276.D
Acq On : 28 Jul 2022 17:45
Operator : JC/MD
Sample : N3953-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

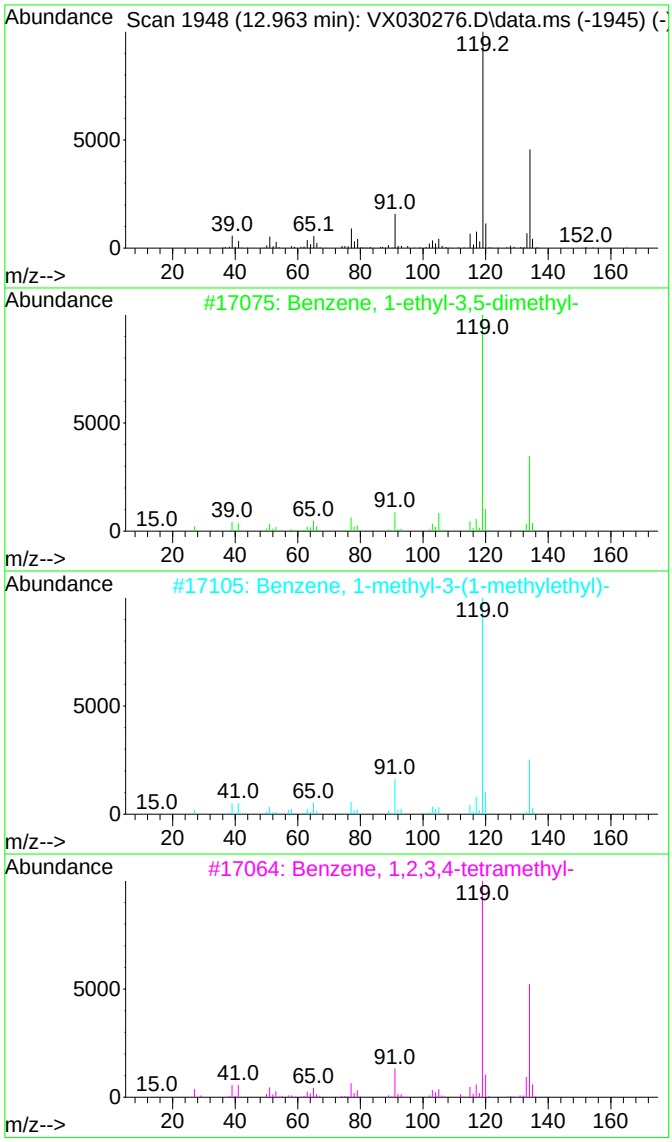
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

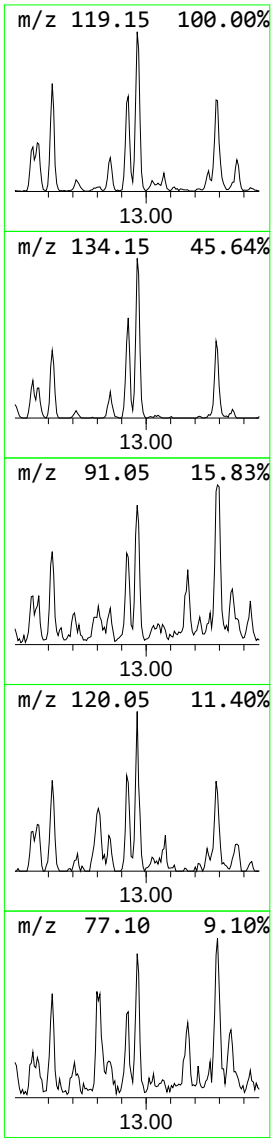
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	9.44 ug/l	123478	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
Data File : VX030276.D
Acq On : 28 Jul 2022 17:45
Operator : JC/MD
Sample : N3953-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

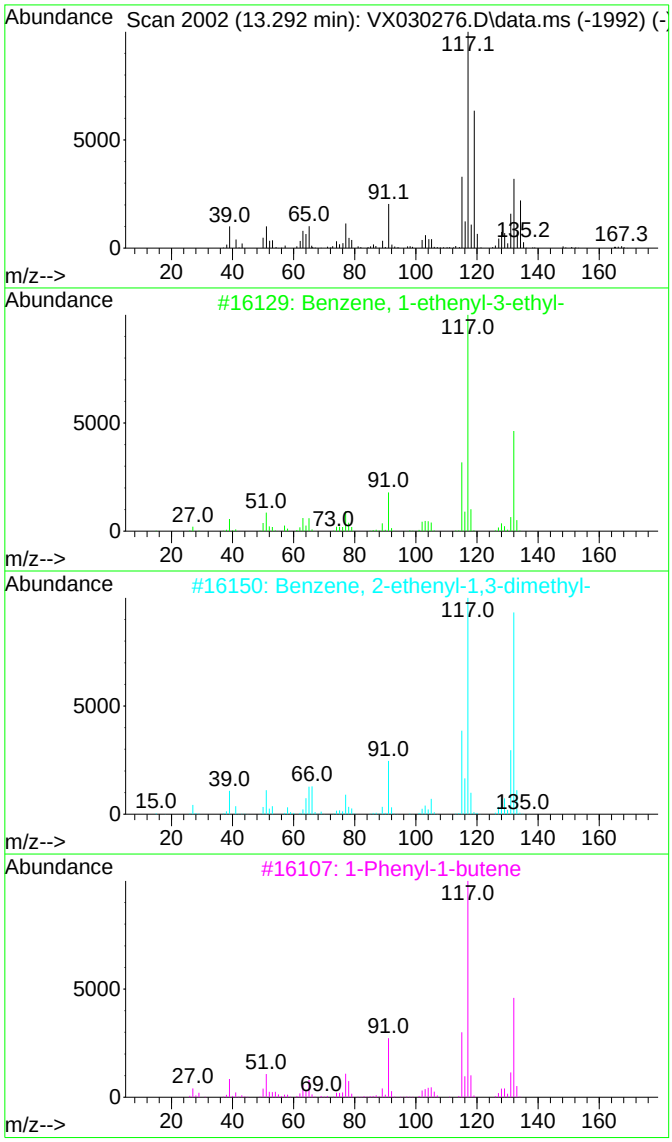
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

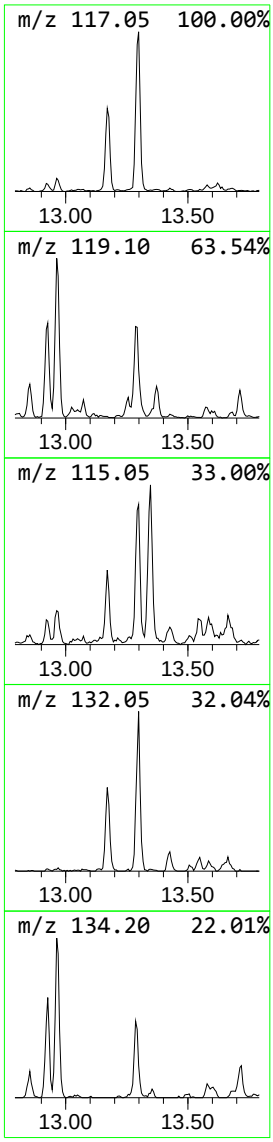
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	20.08 ug/l	262807	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	92
2	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	70
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	70
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	64
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
Data File : VX030276.D
Acq On : 28 Jul 2022 17:45
Operator : JC/MD
Sample : N3953-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

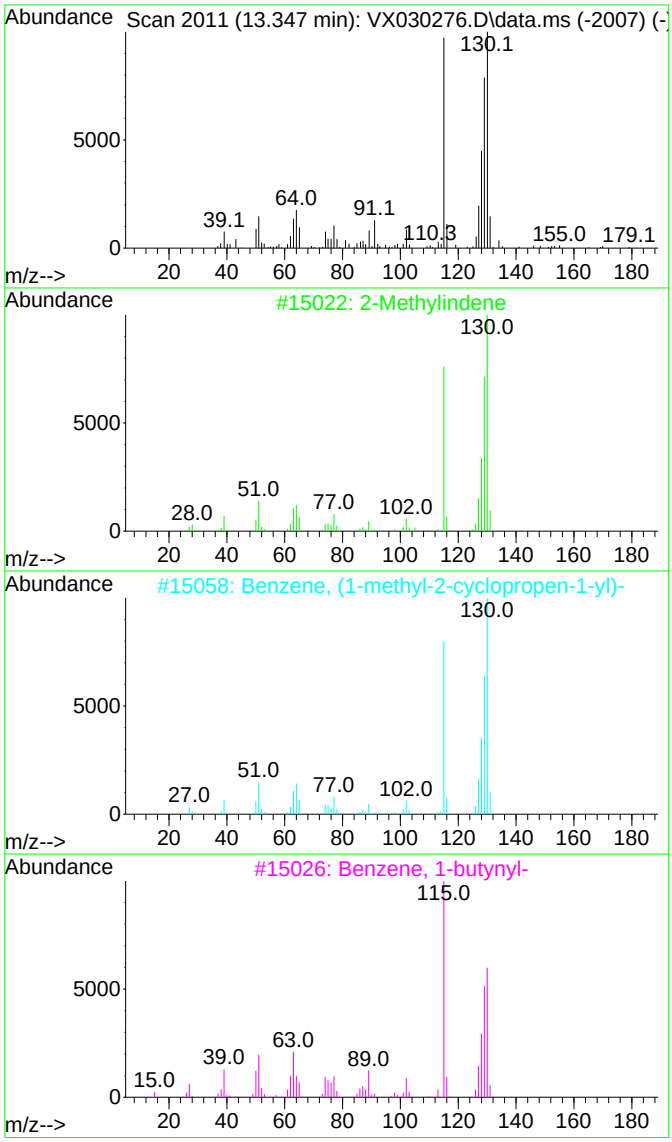
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

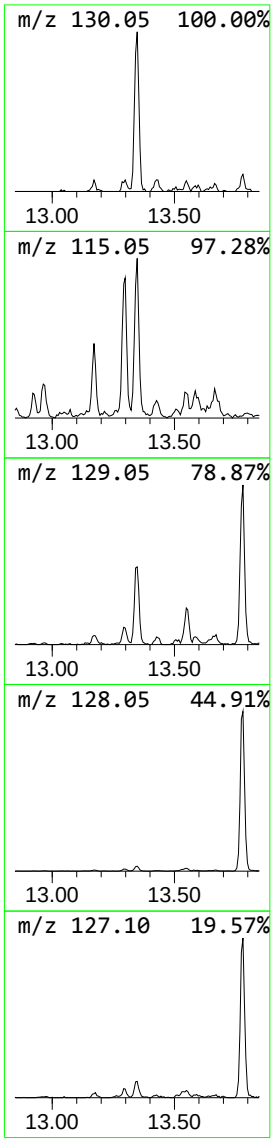
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 2-Methylindene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.347	8.16 ug/l	106712	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylindene	130	C10H10	002177-47-1	96
2	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	96
3	Benzene, 1-butynyl-	130	C10H10	000622-76-4	95
4	Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	93
5	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	93





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
Data File : VX030276.D
Acq On : 28 Jul 2022 17:45
Operator : JC/MD
Sample : N3953-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

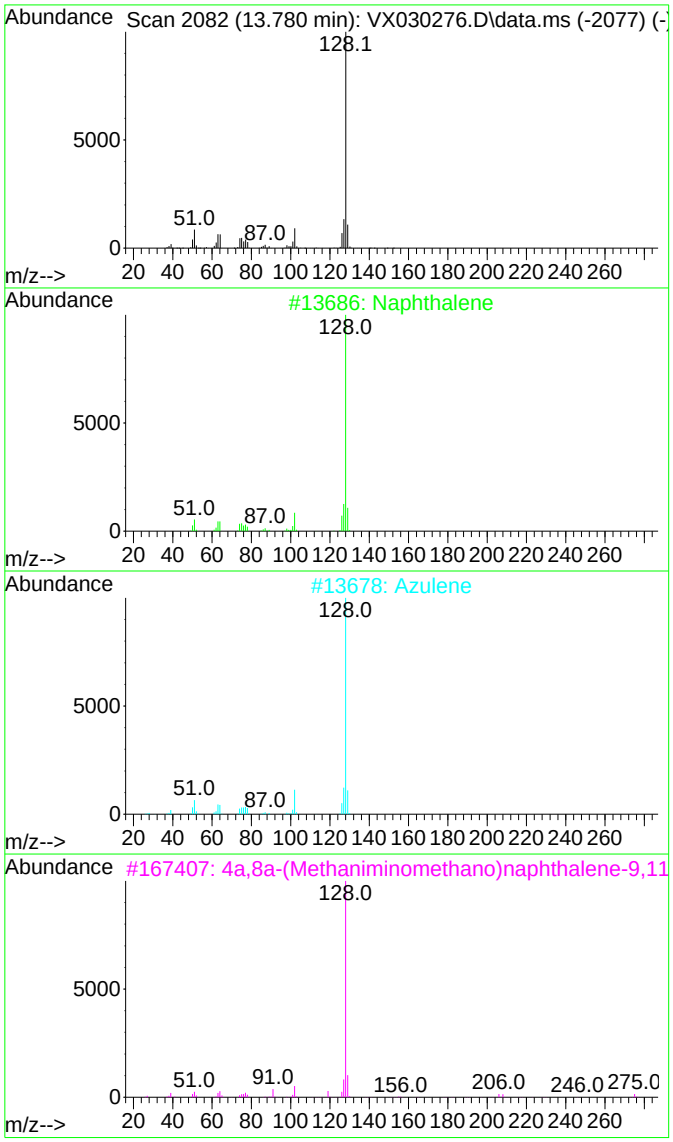
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

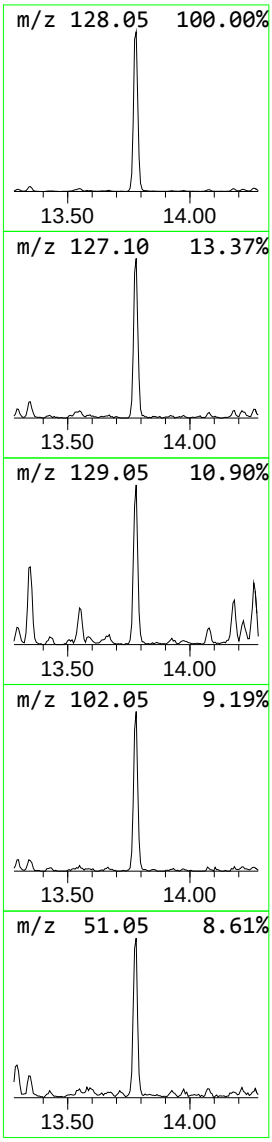
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 4a,8a-(Methaniminomethano)n... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.780	35.45 ug/l	463875	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Azulene	128	C10H8	000275-51-4	94
3	4a,8a-(Methaniminomethano)naphth...	275	C18H13NO2	069915-10-2	64
4	1,3-Dicyanobenzene	128	C8H4N2	000626-17-5	53
5	1H-Indene, 1-methylene-	128	C10H8	002471-84-3	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
Data File : VX030276.D
Acq On : 28 Jul 2022 17:45
Operator : JC/MD
Sample : N3953-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

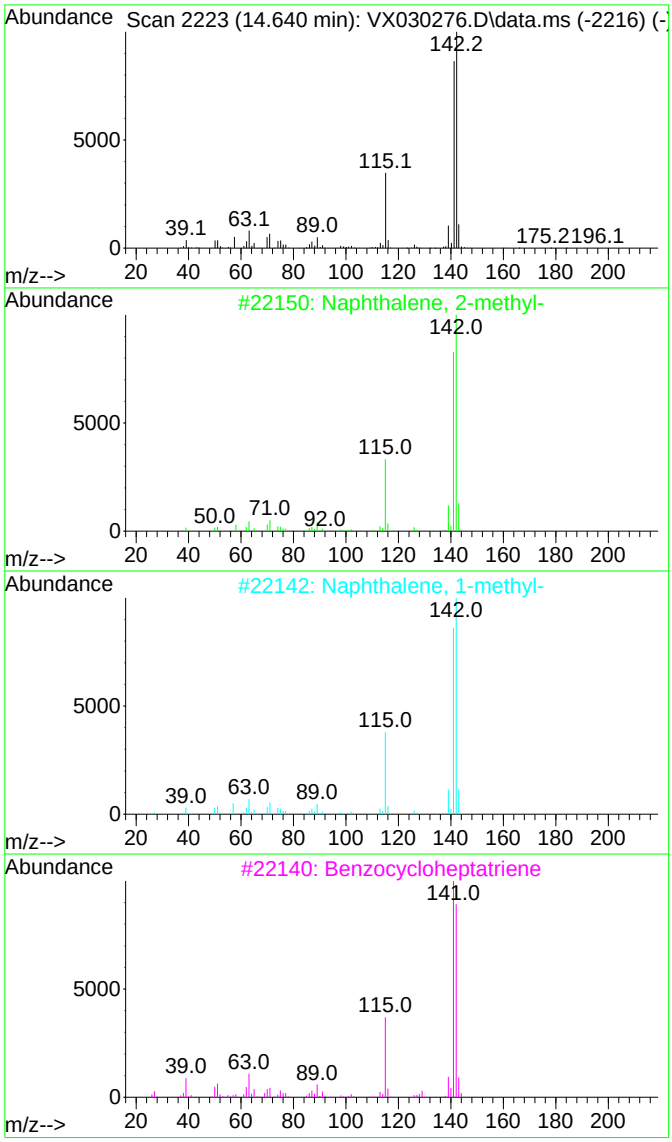
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

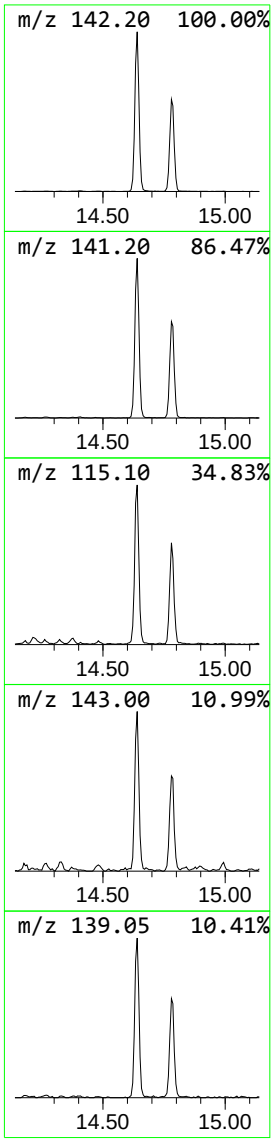
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzocycloheptatriene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.640	48.97 ug/l	640752	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	91
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
Data File : VX030276.D
Acq On : 28 Jul 2022 17:45
Operator : JC/MD
Sample : N3953-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

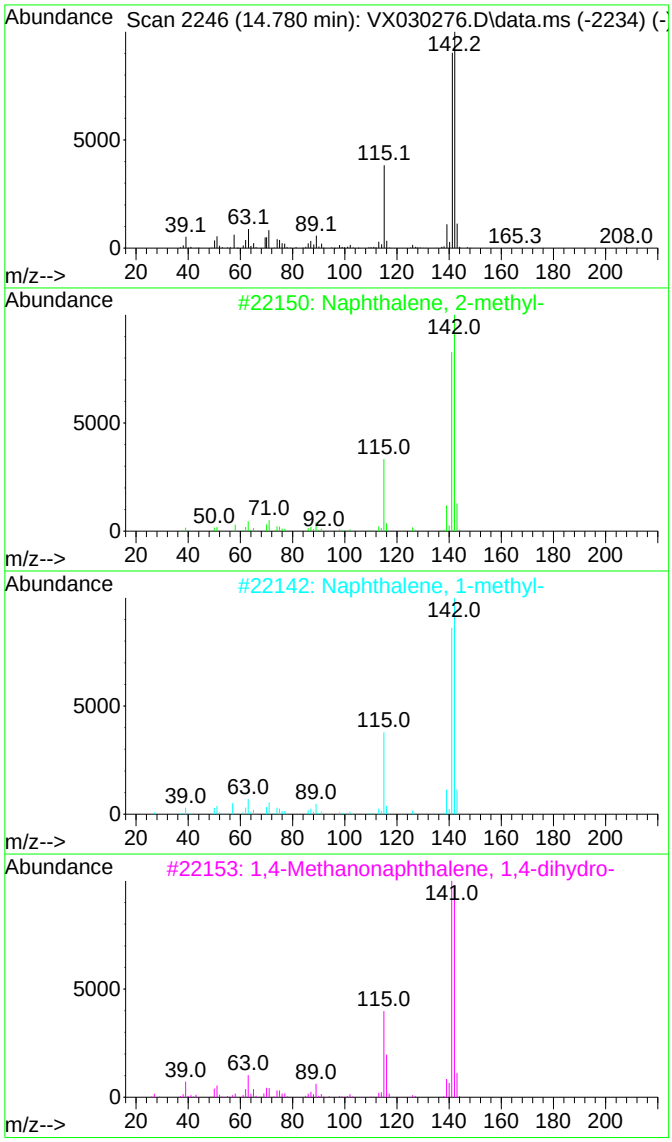
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

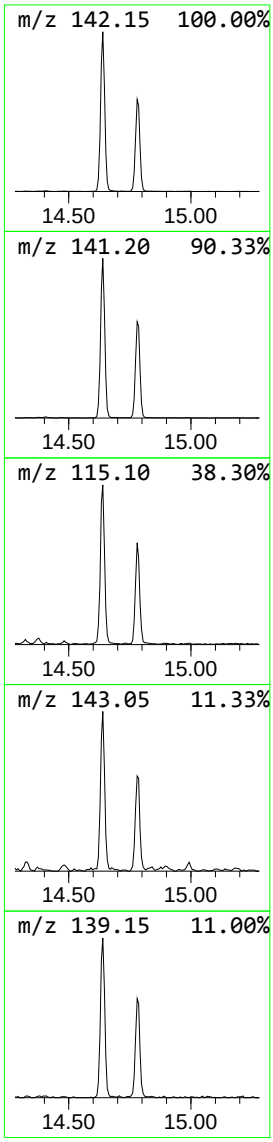
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	35.31 ug/l	462055	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
Data File : VX030276.D
Acq On : 28 Jul 2022 17:45
Operator : JC/MD
Sample : N3953-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IRV-OILY-WATER

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Acetaldehyde	1.453	33.7	ug/l	280493	1	5.562	415791	50.0
Ethanol	2.063	37.7	ug/l	313511	1	5.562	415791	50.0
Pentanal	7.586	12.0	ug/l	140126	2	6.769	582382	50.0
Hexanal	9.525	12.8	ug/l	177736	3	10.055	696853	50.0
Benzene, 1-ethy...	12.963	9.4	ug/l	123478	4	12.024	654250	50.0
Benzene, 1-ethe...	13.292	20.1	ug/l	262807	4	12.024	654250	50.0
2-Methylindene	13.347	8.2	ug/l	106712	4	12.024	654250	50.0
4a,8a-(Methanim...	13.780	35.5	ug/l	463875	4	12.024	654250	50.0
Benzocyclohepta...	14.640	49.0	ug/l	640752	4	12.024	654250	50.0
1,4-Methanonaph...	14.780	35.3	ug/l	462055	4	12.024	654250	50.0